

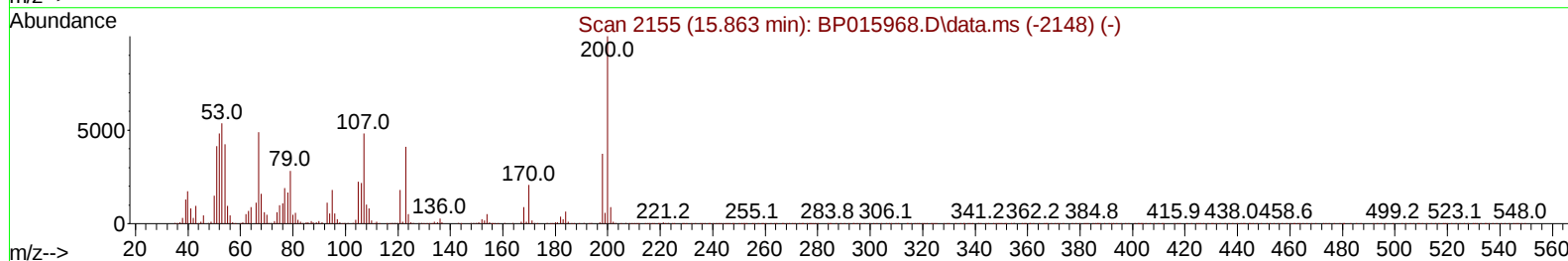
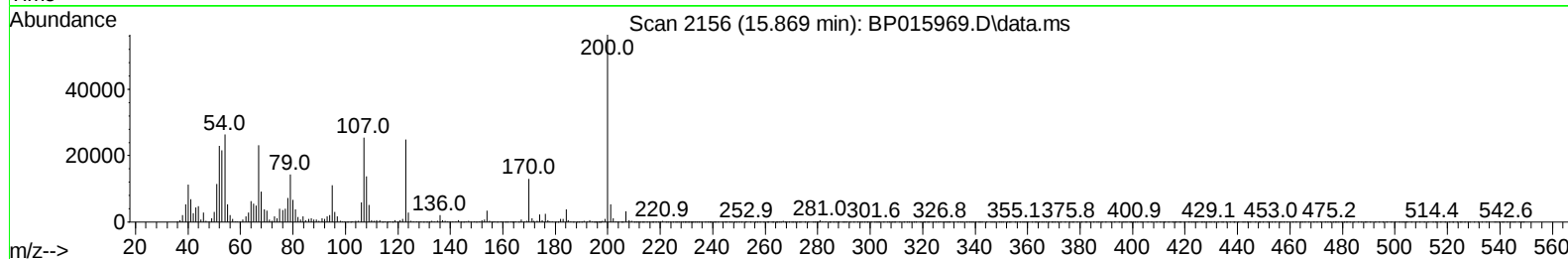
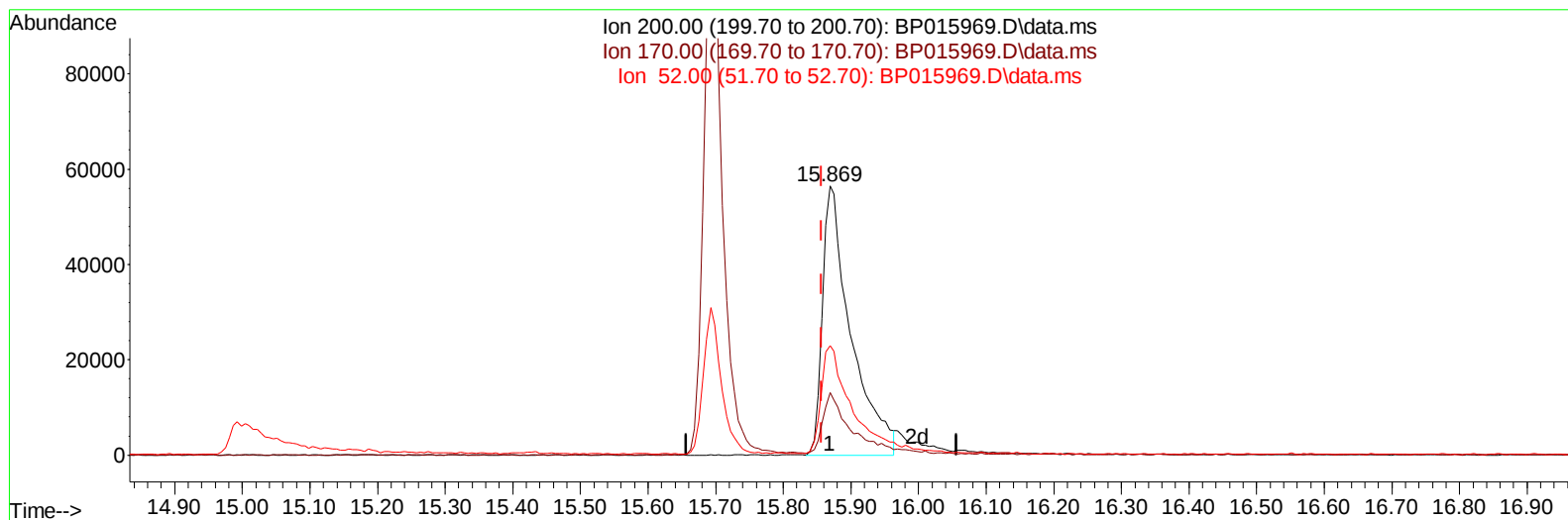
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015969.D
 Acq On : 04 Jul 2023 11:10
 Operator : MA/JU
 Sample : PB153768BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK768

Manual IntegrationsAPPROVED

Quant Time: Jul 04 21:55:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 02 06:12:16 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/05/2023
 Supervised By :mohammad ahmed 07/05/2023



TIC: BP015969.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.869min (+ 0.012) 16.40 ng/u1

response 164475

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	24.30	23.15
52.00	52.20	40.66#
0.00	0.00	0.00

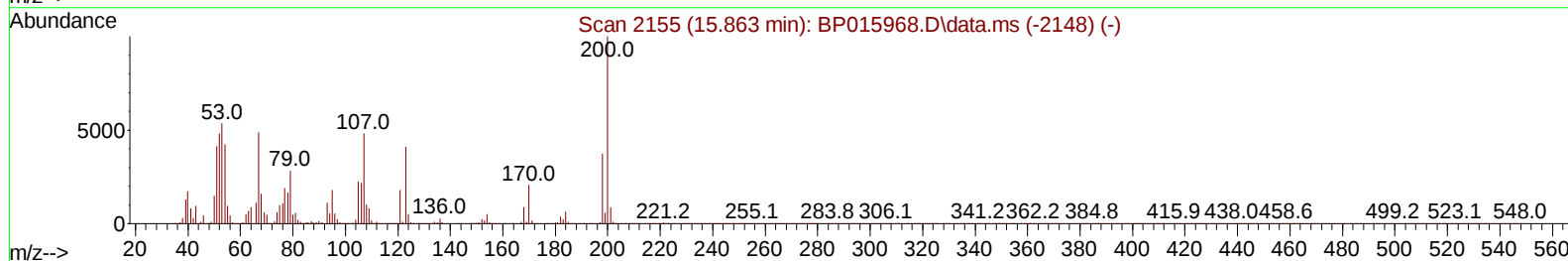
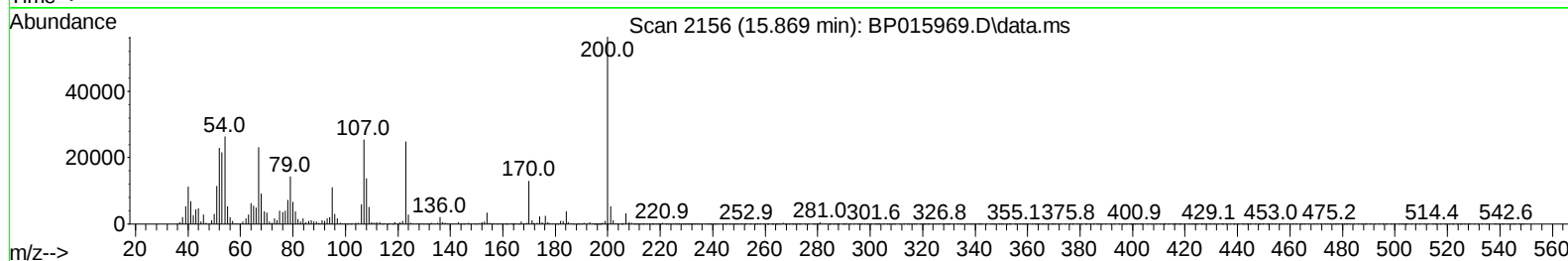
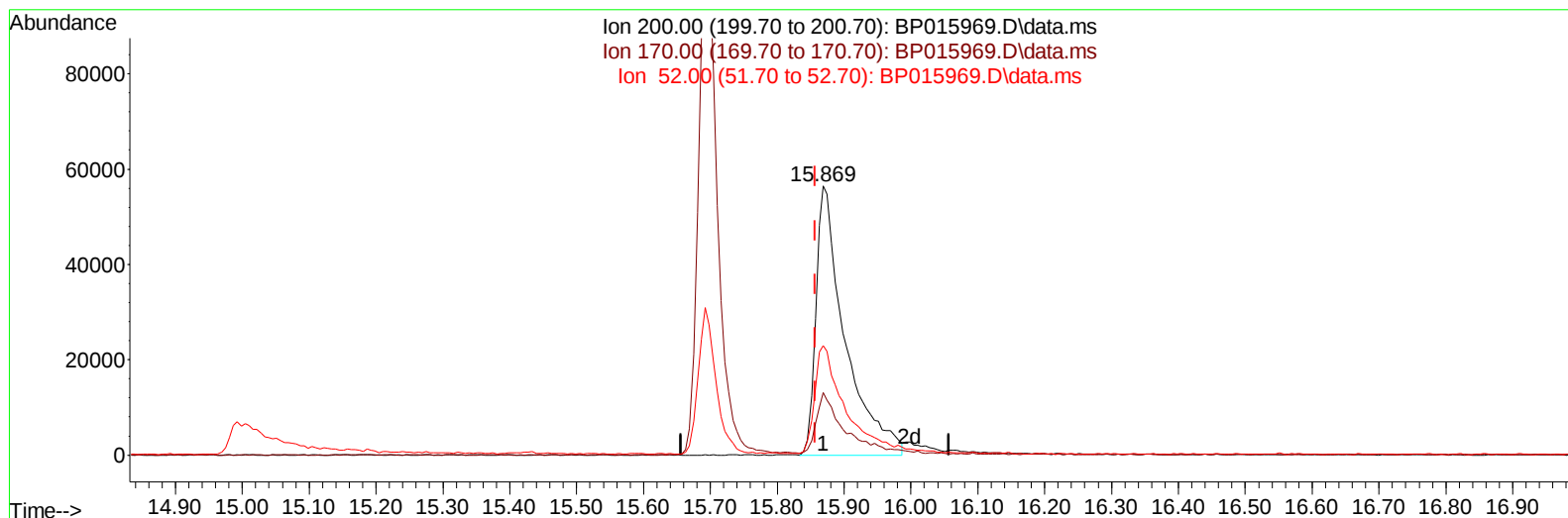
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TIC: BP015969.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.869min (+ 0.012) 16.93 ng/ul m

response 169813

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	24.30	23.15
52.00	52.20	40.66#
0.00	0.00	0.00

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 Operator : MA/JU
 Sample : PB153768BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK768

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/05/2023
 Supervised By :mohammad ahmed 07/05/2023

Quant Time: Jul 04 22:05:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	378156	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	1649575	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.693	164	932089	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.457	188	1630684	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	1232805	20.000	ng/ul	0.00
88) Perylene-d12	24.092	264	1415532	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	53078	5.050	ng/uL	0.00
4) Pyridine-d5	3.811	84	804506	30.364	ng/ul	0.00
7) Phenol-d5	7.228	99	1027337	31.751	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.364	67	698727	34.906	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	817925	33.488	ng/ul	0.00
15) 4-Methylphenol-d8	8.781	113	847215	33.146	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	409820	32.508	ng/ul	0.00
24) 2-Nitrophenol-d4	9.957	143	450529	30.620	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	738904	28.182	ng/ul	0.01
31) 4-Chloroaniline-d4	11.034	131	1073653	31.877	ng/ul	0.00
46) Dimethylphthalate-d6	14.104	166	2321556	32.224	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	2559093	31.599	ng/ul	0.00
54) 4-Nitrophenol-d4	14.993	143	205378	21.602	ng/ul	0.04
60) Fluorene-d10	15.692	176	1796192	30.280	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	169813m	16.933	ng/ul	0.01
73) Anthracene-d10	17.557	188	2444150	32.813	ng/ul	0.00
81) Pyrene-d10	19.786	212	2507717	31.152	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	2371084	33.206	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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